

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111022\
 Data File : BM037451.D
 Acq On : 10 Nov 2022 14:57
 Operator : CG/JU
 Sample : N5436-03DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 250FERRYBLVD-TFS-03DL

Quant Time: Nov 10 17:21:33 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 09 02:08:14 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 11/11/2022
 Supervised By :Jagrut Upadhyay 11/11/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	244439	20.000	ng	-0.01
21) Naphthalene-d8	10.610	136	952372	20.000	ng	-0.02
39) Acenaphthene-d10	14.451	164	546314	20.000	ng	-0.01
64) Phenanthrene-d10	17.192	188	1169257	20.000	ng	-0.01
76) Chrysene-d12	21.374	240	1252458	20.000	ng	0.00
86) Perylene-d12	23.703	264	1338787	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.387	112	274335	19.389	ng	0.00
7) Phenol-d6	6.992	99	327088	17.033	ng	-0.01
23) Nitrobenzene-d5	8.981	82	199276	10.557	ng	-0.02
42) 2,4,6-Tribromophenol	15.939	330	136198	21.155	ng	-0.02
45) 2-Fluorobiphenyl	13.074	172	458928	11.162	ng	-0.02
79) Terphenyl-d14	19.815	244	764399	10.945	ng	-0.01
Target Compounds						Qvalue
20) 3+4-Methylphenols	8.598	107	50770	2.629	ng	96
27) 2,4-Dimethylphenol	9.798	122	67612	5.318	ng	95
31) Naphthalene	10.663	128	155371	2.880	ng	99
49) Acenaphthylene	14.174	152	120607	2.215	ng	99
71) Phenanthrene	17.233	178	912556	12.580	ng	100
72) Anthracene	17.327	178	268767	3.914	ng	99
75) Fluoranthene	19.251	202	3482821	45.419	ng	99
78) Pyrene	19.615	202	3514330	36.490	ng	98
81) Benzo(a)anthracene	21.356	228	1404144	15.747	ng	99
83) Chrysene	21.409	228	1286329	14.974	ng	98
87) Indeno(1,2,3-cd)pyrene	26.109	276	550040	5.018	ng	95
88) Benzo(b)fluoranthene	22.997	252	1269444	13.658	ng	98
89) Benzo(k)fluoranthene	23.039	252	459513m	4.847	ng	
90) Benzo(a)pyrene	23.603	252	910711	12.026	ng	99
92) Benzo(g,h,i)perylene	26.856	276	521779	5.889	ng	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111022\
Data File : BM037451.D
Acq On : 10 Nov 2022 14:57
Operator : CG/JU
Sample : N5436-03DL 5X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
250FERRYBLVD-TFS-03DL

Quant Time: Nov 10 17:21:33 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 09 02:08:14 2022
Response via : Initial Calibration

Manual Integrations
APPROVED

Reviewed By : Christian Giraldo 11/11/2022
Supervised By : Jagrut Upadhyay 11/11/2022

