

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102623\
 Data File : BM042468.D
 Acq On : 27 Oct 2023 00:42
 Operator : MA/JU
 Sample : 04805-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DUP-02

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 10/27/2023
 Supervised By :mohammad ahmed 10/27/2023

Quant Time: Oct 27 01:13:05 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 21 02:32:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.992	152	211450	20.000	ng	0.00
21) Naphthalene-d8	10.810	136	837873	20.000	ng	-0.01
39) Acenaphthene-d10	14.633	164	488086	20.000	ng	-0.01
64) Phenanthrene-d10	17.386	188	932526	20.000	ng	-0.01
76) Chrysene-d12	21.562	240	621741	20.000	ng	-0.02
86) Perylene-d12	24.027	264	657640	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	875939	58.547	ng	0.00
7) Phenol-d6	7.145	99	1164198	58.065	ng	0.00
23) Nitrobenzene-d5	9.186	82	737225	38.908	ng	-0.01
42) 2,4,6-Tribromophenol	16.127	330	360849	71.428	ng	-0.01
45) 2-Fluorobiphenyl	13.257	172	1456245	40.177	ng	-0.01
79) Terphenyl-d14	19.998	244	1638696	46.743	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	10.863	128	125730	2.921	ng	99
38) 1-Methylnaphthalene	12.686	142	104261	3.842	ng	100
49) Acenaphthylene	14.363	152	306219	6.677	ng	98
52) Acenaphthene	14.698	154	58860	2.123	ng	99
55) Dibenzofuran	15.033	168	92494	2.109	ng	98
58) Fluorene	15.680	166	104306	3.014	ng	99
71) Phenanthrene	17.427	178	1217329	23.996	ng	99
72) Anthracene	17.521	178	500227	9.956	ng	99
75) Fluoranthene	19.439	202	1541621	27.593	ng	98
78) Pyrene	19.803	202	1699215	39.110	ng	100
81) Benzo(a)anthracene	21.550	228	749428	17.980	ng	95
83) Chrysene	21.603	228	676887m	16.909	ng	
87) Indeno(1,2,3-cd)pyrene	26.597	276	316450	6.584	ng	# 92
88) Benzo(b)fluoranthene	23.268	252	655409m	16.557	ng	
89) Benzo(k)fluoranthene	23.309	252	251334m	6.115	ng	
90) Benzo(a)pyrene	23.915	252	602522	15.814	ng	98
91) Dibenzo(a,h)anthracene	26.615	278	92374	2.312	ng	# 88
92) Benzo(g,h,i)perylene	27.409	276	315075	8.119	ng	92

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

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