

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030420\
 Data File : BF119191.D
 Acq On : 4 Mar 2020 13:27
 Operator : CG/JU
 Sample : L1762-04
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-05-030220

Manual Integrations
 APPROVED

mohammad
 3/5/2020 3:54:16 PM

Quant Time: Mar 04 14:46:10 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 26 16:03:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	145771	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	552705	20.00	ng	-0.01
39) Acenaphthene-d10	9.77	164	289237	20.00	ng	-0.01
64) Phenanthrene-d10	11.26	188	477064	20.00	ng	0.00
76) Chrysene-d12	13.89	240	268011	20.00	ng	-0.01
87) Perylene-d12	15.32	264	309074	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	834615	95.68	ng	0.00
7) Phenol-d6	6.39	99	1020310	96.18	ng	-0.01
23) Nitrobenzene-d5	7.30	82	678676	64.83	ng	-0.02
42) 2,4,6-Tribromophenol	10.56	330	327547	86.57	ng	-0.01
45) 2-Fluorobiphenyl	9.10	172	1351450	68.83	ng	-0.01
79) Terphenyl-d14	12.84	244	1269680	81.56	ng	-0.01
Target Compounds						Qvalue
50) Dimethylphthalate	9.49	163	215908	9.762	ng	99
71) Phenanthrene	11.28	178	91083	3.501	ng	96
75) Fluoranthene	12.47	202	145660	5.274	ng	98
78) Pyrene	12.70	202	139383	6.477	ng	97
81) Benzo(a)anthracene	13.89	228	54333	2.975	ng	97
83) Chrysene	13.92	228	54508	2.996	ng	97
88) Benzo(b)fluoranthene	14.92	252	62955m	3.090	ng	
90) Benzo(a)pyrene	15.27	252	49289	2.681	ng	98

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

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