

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120518\
 Data File : BF111115.D
 Acq On : 5 Dec 2018 18:58
 Operator : JU/SJ
 Sample : J5764-21 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-113018-A

Manual Integrations
 APPROVED

mohammad

12/6/2018 12:11:50 PM

Quant Time: Dec 06 01:18:17 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 05 17:30:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	578502	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	2258448	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	1008775	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	1543379	20.00	ng	0.00
76) Chrysene-d12	13.96	240	1092094	20.00	ng	0.00
87) Perylene-d12	15.39	264	807845	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	1759246	47.45	ng	0.00
7) Phenol-d6	6.43	99	2300092	48.16	ng	0.00
23) Nitrobenzene-d5	7.36	82	1385070	36.31	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	366793	45.74	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	2389351	35.47	ng	0.00
79) Terphenyl-d14	12.91	244	1683022	31.56	ng	0.00
Target Compounds						
10) Phenol	6.43	94	137825	2.647	ng	90
50) Dimethylphthalate	9.55	163	163723	2.332	ng	99
71) Phenanthrene	11.35	178	338558	4.560	ng	93
75) Fluoranthene	12.53	202	356988	4.914	ng	91
78) Pyrene	12.76	202	332690	3.910	ng	96
81) Benzo(a)anthracene	13.94	228	141768	2.216	ng	98
83) Chrysene	13.98	228	140219	2.207	ng	96
88) Benzo(b)fluoranthene	14.97	252	141338m	3.088	ng	
90) Benzo(a)pyrene	15.33	252	94194	2.243	ng	# 93

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

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