

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112718\
 Data File : BF111004.D
 Acq On : 27 Nov 2018 12:43
 Operator : JU/SJ
 Sample : J6028-08RE
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-3RE

Manual Integrations
 APPROVED

mohammad
 11/28/2018 4:42:09 PM

Quant Time: Nov 27 14:41:54 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 27 06:43:50 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	45475	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	162945	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	65035	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	126412	20.00	ng	0.00
76) Chrysene-d12	14.14	240	109178	20.00	ng	0.00
87) Perylene-d12	15.67	264	73458	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	289042	105.40	ng	0.01
7) Phenol-d6	6.57	99	342017	95.71	ng	0.00
23) Nitrobenzene-d5	7.50	82	210626	73.98	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	68709	106.38	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	335925	74.83	ng	0.00
79) Terphenyl-d14	13.06	244	346897	61.85	ng	0.00
Target Compounds						
10) Phenol	6.58	94	17959	4.462	ng	87
50) Dimethylphthalate	9.69	163	45430	9.525	ng	99
71) Phenanthrene	11.50	178	68885	10.289	ng	99
72) Anthracene	11.56	178	22894	3.358	ng	98
75) Fluoranthene	12.70	202	144582	21.011	ng	99
78) Pyrene	12.93	202	135013	16.932	ng	98
81) Benzo(a)anthracene	14.13	228	64519	10.010	ng	99
83) Chrysene	14.16	228	49899	7.864	ng	96
86) Indeno(1,2,3-cd)pyrene	17.28	276	20012	4.394	ng	# 85
88) Benzo(b)fluoranthene	15.22	252	51372m	11.904	ng	
89) Benzo(k)fluoranthene	15.24	252	16613m	3.740	ng	
90) Benzo(a)pyrene	15.62	252	32435	8.075	ng	# 96
92) Benzo(a,h,i)perylene	17.77	276	21189	6.631	ng	# 89

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

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