

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112618\
 Data File : BF110968.D
 Acq On : 26 Nov 2018 17:20
 Operator : JU/SJ
 Sample : J6030-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DTW-03

Manual Integrations
 APPROVED

mohammad
 11/27/2018 11:47:39 AM

Quant Time: Nov 27 10:41:24 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 26 15:26:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	62765	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	265749	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	131337	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	216151	20.00	ng	0.00
76) Chrysene-d12	14.13	240	131664	20.00	ng	0.00
87) Perylene-d12	15.67	264	132115	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	304729	80.51	ng	0.01
7) Phenol-d6	6.59	99	263900	53.50	ng	0.02
23) Nitrobenzene-d5	7.51	82	354594	76.36	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	141219	108.27	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	658745	72.67	ng	0.00
79) Terphenyl-d14	13.06	244	533662	78.90	ng	0.00
Target Compounds						
6) Aniline	6.62	93	40315	6.612	ng	80
10) Phenol	6.60	94	48990	8.820	ng	83
17) 2-Methylphenol	7.20	107	20399	5.720	ng	80
20) 3+4-Methylphenols	7.36	107	100587m	21.217	ng	
22) Acetophenone	7.35	105	231779	37.209	ng	93
27) 2,4-Dimethylphenol	7.88	122	32266	9.107	ng	99
31) Naphthalene	8.26	128	3001696	240.963	ng	98
32) Benzoic acid	8.02	122	118054m	44.818	ng	
37) 2-Methylnaphthalene	8.93	142	101876	12.382	ng	99
38) 1-Methylnaphthalene	9.03	142	97865	12.228	ng	95
46) 1,1'-Biphenyl	9.40	154	36393	3.006	ng	100
55) Dibenzofuran	10.19	168	82801	7.092	ng	97
58) Fluorene	10.53	166	67326	7.412	ng	99
71) Phenanthrene	11.50	178	213090	18.614	ng	99

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

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