

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092519\
 Data File : BF117262.D
 Acq On : 26 Sep 2019 2:37
 Operator : JU
 Sample : K5008-03DL 2X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-3DL

Manual Integrations
 APPROVED

mohammad
 9/26/2019 3:54:05 PM

Quant Time: Sep 26 10:59:13 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 13 16:23:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	37631	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	141596	20.00	ng	0.00
39) Acenaphthene-d10	9.86	164	66094	20.00	ng	0.00
64) Phenanthrene-d10	11.35	188	84862	20.00	ng	0.02
76) Chrysene-d12	13.98	240	66919	20.00	ng	0.00
87) Perylene-d12	15.43	264	55441	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	109556	45.45	ng	0.00
7) Phenol-d6	6.45	99	148897	47.80	ng	0.00
23) Nitrobenzene-d5	7.38	82	91033	33.78	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	21560	39.03	ng	0.01
45) 2-Fluorobiphenyl	9.17	172	146342	34.62	ng	0.00
79) Terphenyl-d14	12.92	244	104975	29.32	ng	0.00
Target Compounds						
10) Phenol	6.47	94	7526	2.192	ng	79
31) Naphthalene	8.12	128	199909	29.356	ng	# 94
37) 2-Methylnaphthalene	8.83	142	722201	157.489	ng	99
38) 1-Methylnaphthalene	8.93	142	600599m	139.840	ng	
46) 1,1'-Biphenyl	9.28	154	63311	12.212	ng	99
49) Acenaphthylene	9.73	152	20455m	3.337	ng	
50) Dimethylphthalate	9.57	163	24466	5.228	ng	# 64
52) Acenaphthene	9.90	154	41350	11.380	ng	# 73
55) Dibenzofuran	10.07	168	56053	10.456	ng	# 5
58) Fluorene	10.42	166	92549	21.431	ng	# 92
71) Phenanthrene	11.37	178	256234	53.159	ng	# 96
72) Anthracene	11.43	178	19799	4.023	ng	# 60
78) Pvrene	12.79	202	16984	3.025	ng	# 94

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

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