

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP092419\
 Data File : BP000391.D
 Acq On : 24 Sep 2019 18:12
 Operator : JU
 Sample : K5008-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-1

Manual Integrations
 APPROVED

mohammad
 9/26/2019 8:24:53 AM

Quant Time: Sep 25 11:39:05 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP091619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 19 14:44:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	252510	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	688171	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	387804	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	577980	20.00	ng	0.00
76) Chrysene-d12	13.99	240	710004	20.00	ng	0.00
87) Perylene-d12	15.47	264	753885	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	1605481	109.76	ng	0.00
7) Phenol-d6	6.43	99	2006803	111.93	ng	0.00
23) Nitrobenzene-d5	7.35	82	1100336	103.16	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	455362	118.38	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	1762136	81.78	ng	0.00
79) Terphenyl-d14	12.93	244	2027843	59.92	ng	0.00
Target Compounds						
10) Phenol	6.44	94	120041	5.893	ng	98
31) Naphthalene	8.10	128	1982189	62.193	ng	96
37) 2-Methylnaphthalene	8.80	142	3139536	144.098	ng	96
38) 1-Methylnaphthalene	8.90	142	2054400m	100.660	ng	
46) 1,1'-Biphenyl	9.26	154	473228	17.640	ng	98
49) Acenaphthylene	9.70	152	109714	3.243	ng	# 1
50) Dimethylphthalate	9.55	163	252201	9.545	ng	# 81
52) Acenaphthene	9.87	154	183857	8.613	ng	# 64
55) Dibenzofuran	10.05	168	200001m	6.627	ng	
58) Fluorene	10.39	166	442780	18.996	ng	# 92
71) Phenanthrene	11.36	178	1029111	35.531	ng	# 96
72) Anthracene	11.41	178	76619	2.570	ng	# 67
73) Carbazole	11.56	167	59277	2.208	ng	# 24
75) Fluoranthene	12.56	202	117482	3.777	ng	91
78) Pyrene	12.79	202	223868	4.214	ng	96

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

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