

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP100419\
 Data File : BP000511.D
 Acq On : 04 Oct 2019 16:48
 Operator : JU
 Sample : K5095-05
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COMPOSITE-5

Quant Time: Oct 05 00:02:22 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 01 18:03:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	190542	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	618776	20.00	ng	0.00
39) Acenaphthene-d10	9.81	164	413917	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	708285	20.00	ng	0.00
76) Chrysene-d12	13.98	240	410128	20.00	ng	0.00
87) Perylene-d12	15.46	264	482475	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	996802	84.09	ng	0.00
7) Phenol-d6	6.41	99	819680	57.38	ng	0.00
23) Nitrobenzene-d5	7.33	82	721287	71.19	ng	0.00
42) 2,4,6-Tribromophenol	10.61	330	600217	136.08	ng	0.00
45) 2-Fluorobiphenyl	9.13	172	1855449	72.53	ng	0.00
79) Terphenyl-d14	12.92	244	1772328	87.50	ng	0.00
Target Compounds						
10) Phenol	6.42	94	173061	11.833	ng	Qvalue 89
13) 1,4-Dichlorobenzene	6.79	146	45672	3.201	ng	97
14) 1,2-Dichlorobenzene	6.94	146	86234	6.545	ng	99
15) Benzyl Alcohol	6.90	79	41597	4.509	ng	97
17) 2-Methylphenol	7.02	107	69897	7.230	ng	# 86
20) 3+4-Methylphenols	7.17	107	157561	13.927	ng	# 60
30) 1,2,4-Trichlorobenzene	7.99	180	69541	6.648	ng	99
31) Naphthalene	8.07	128	381529	13.202	ng	100
37) 2-Methylnaphthalene	8.76	142	87361	4.503	ng	98
38) 1-Methylnaphthalene	8.86	142	51752	2.823	ng	99
50) Dimethylphthalate	9.53	163	383964	13.350	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.14	232	19023	2.703	ng	# 92
70) Pentachlorophenol	11.12	266	24353	6.692	ng	100
74) Di-n-butylphthalate	11.89	149	106241	2.667	ng	99

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

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