

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081617\
 Data File : BF097763.D
 Acq On : 17 Aug 2017 3:07
 Operator : SJ/JU
 Sample : I4773-04
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-J10-SSW

Manual Integrations
 APPROVED

Sohil
 8/17/2017 5:58:04 PM

Quant Time: Aug 17 07:51:42 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 15 13:25:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	127435	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	502521	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	205688	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	303729	20.00	ng	0.00
75) Chrysene-d12	13.93	240	245794	20.00	ng	0.00
86) Perylene-d12	15.36	264	214796	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	1033286	123.33	ng	0.00
7) Phenol-d6	6.42	99	1366023	120.00	ng	0.00
23) Nitrobenzene-d5	7.35	82	851256	92.02	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	206800	115.88	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	1133628	80.97	ng	0.00
78) Terphenyl-d14	12.89	244	519347	44.06	ng	0.00
Target Compounds						
10) Phenol	6.42	94	46726	3.510	ng	91
12) 1,3-Dichlorobenzene	6.72	146	37594	3.956	ng	# 92
13) 1,4-Dichlorobenzene	6.80	146	121602m	12.581	ng	
14) 1,2-Dichlorobenzene	6.95	146	493154	55.333	ng	98
30) 1,2,4-Trichlorobenzene	8.00	180	18626	2.523	ng	97
31) Naphthalene	8.09	128	869178	33.317	ng	99
32) Benzoic acid	7.79	122	19222m	9.254	ng	
37) 2-Methylnaphthalene	8.77	142	103934	6.498	ng	97
49) Dimethylphthalate	9.53	163	287503	19.122	ng	100
70) Phenanthrene	11.32	178	33147	2.048	ng	98

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

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