

Data Path : Z:\HPCHEM1\BNA F\DATA\BF072517\
 Data File : BF097060.D
 Acq On : 26 Jul 2017 6:34
 Operator : SJ/JU
 Sample : I4370-06
 Misc :
 ALS Vial : 34 Sample Multiplier: 2

Instrument :
 BNA_F
 ClientSampleId :
 E2-N5-NWSW

Manual Integrations
 APPROVED

Sohil
 7/26/2017 5:58:37 PM

Quant Time: Jul 26 15:37:14 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 25 14:11:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.50	152	207841	40.00	ng	0.00
21) Naphthalene-d8	7.78	136	811864	40.00	ng	0.00
38) Acenaphthene-d10	9.53	164	312553	40.00	ng	0.00
63) Phenanthrene-d10	11.00	188	445731	40.00	ng	0.00
75) Chrysene-d12	13.63	240	358091	40.00	ng	0.00
86) Perylene-d12	14.97	264	271103	40.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.11	112	841274	122.04	ng	0.01
7) Phenol-d6	6.17	99	1082072	137.50	ng	0.00
23) Nitrobenzene-d5	7.07	82	582854	84.23	ng	0.00
41) 2,4,6-Tribromophenol	10.32	330	142398	115.33	ng	0.00
44) 2-Fluorobiphenyl	8.86	172	816063	88.62	ng	0.00
78) Terphenyl-d14	12.59	244	477231	54.34	ng	0.00
Target Compounds						
10) Phenol	6.18	94	26770	2.868	ng	96
11) bis(2-Chloroethyl)ether	6.24	93	55795	7.557	ng	89
12) 1,3-Dichlorobenzene	6.44	146	34305	4.319	ng	98
13) 1,4-Dichlorobenzene	6.52	146	59782	7.594	ng	96
14) 1,2-Dichlorobenzene	6.67	146	118385	17.223	ng	97
16) 2,2'-oxybis(1-Chloropropan	6.79	45	54678	3.692	ng	92
49) Dimethylphthalate	9.26	163	100171	8.441	ng	100
70) Phenanthrene	11.02	178	43852	3.672	ng	96
74) Fluoranthene	12.20	202	70946	6.065	ng	97
77) Pyrene	12.43	202	71394	4.871	ng	97
80) Benzo(a)anthracene	13.62	228	34559	2.983	ng	97
82) Chrysene	13.64	228	33752m	2.984	ng	
87) Benzo(b)fluoranthene	14.62	252	36731m	4.508	ng	
89) Benzo(a)pyrene	14.92	252	26464	3.549	ng	# 94

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

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