

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE090517\
 Data File : BE093949.D
 Acq On : 5 Sep 2017 22:43
 Operator : SJ/JU
 Sample : I5066-05
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 TW-1

Manual Integrations
 APPROVED

Sohil
 9/7/2017 6:32:01 PM

Quant Time: Sep 06 06:58:26 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE090517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 05 13:34:56 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2090	0.40	ng	0.00
7) Naphthalene-d8	10.71	136	10228	0.40	ng	0.00
13) Acenaphthene-d10	14.52	164	6136	0.40	ng	0.00
19) Phenanthrene-d10	17.23	188	10673	0.40	ng	-0.03
27) Chrysene-d12	21.41	240	14636	0.40	ng	-0.01
34) Perylene-d12	23.93	264	12883	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.53	112	1491	0.22	ng	0.00
5) Phenol-d6	7.12	99	1379	0.14	ng	0.00
8) Nitrobenzene-d5	9.08	82	2951	0.34	ng	0.00
11) 2-Methylnaphthalene-d10	12.29	152	5304	0.32	ng	0.00
14) 2,4,6-Tribromophenol	16.03	330	618	0.33	ng	0.00
15) 2-Fluorobiphenyl	13.16	172	7726	0.32	ng	0.00
25) Fluoranthene-d10	19.28	212	57764	0.36	ng	0.00
29) Terphenyl-d14	19.86	244	11679	0.43	ng	-0.01
Target Compounds						
16) Acenaphthylene	14.24	152	1069	0.035	ng	Qvalue # 84
24) Anthracene	17.37	178	1009m	0.031	ng	
26) Fluoranthene	19.30	202	5292	0.111	ng	99
28) Pyrene	19.67	202	5190	0.103	ng	98
30) Benzo(a)anthracene	21.40	228	3847	0.086	ng	# 77
31) Chrysene	21.45	228	3464	0.073	ng	# 80
32) Bis(2-ethylhexyl)phthalate	21.30	149	13451	0.123	ng	# 100
33) Indeno(1,2,3-cd)pyrene	26.59	276	3540	0.112	ng	92
35) Benzo(b)fluoranthene	23.16	252	6318	0.154	ng	# 74
36) Benzo(k)fluoranthene	23.20	252	2083m	0.041	ng	
37) Benzo(a)pyrene	23.82	252	5096	0.122	ng	# 74
39) Benzo(a,h,i)perylene	27.41	276	3538	0.121	ng	# 78

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

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