

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE121117\
 Data File : BE094973.D
 Acq On : 11 Dec 2017 14:38
 Operator : SJ/JU
 Sample : I6740-12DL 2X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 FT-PZ-460I-20171204DL

Manual Integrations
 APPROVED

Sohil
 12/11/2017 6:24:10 PM

Quant Time: Dec 11 16:01:20 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE113017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 01 13:33:28 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.48	152	8503	0.40	ng	0.00
7) Naphthalene-d8	11.33	136	17708	0.40	ng	0.00
13) Acenaphthene-d10	15.08	164	9687	0.40	ng	0.00
19) Phenanthrene-d10	17.79	188	24723	0.40	ng	0.02
27) Chrysene-d12	21.89	240	26029	0.40	ng	0.00
34) Perylene-d12	24.57	264	21365	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.95	112	883	0.07	ng	0.00
5) Phenol-d6	7.60	99	975	0.05	ng	0.00
8) Nitrobenzene-d5	9.66	82	2605	0.25	ng	0.00
11) 2-Methylnaphthalene-d10	12.87	152	4649	0.16	ng	0.00
14) 2,4,6-Tribromophenol	16.55	330	390	0.26	ng	0.02
15) 2-Fluorobiphenyl	13.71	172	12148	0.29	ng	0.00
25) Fluoranthene-d10	19.77	212	51867	0.15	ng	0.00
29) Terphenyl-d14	20.33	244	14674	0.31	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.66	88	14446	1.624	ng	Qvalue # 87
22) Pentachlorophenol	17.43	266	141	0.165	ng	91
23) Phenanthrene	17.82	178	2604	0.033	ng	# 93
30) Benzo(a)anthracene	21.87	228	2154	0.027	ng	# 66
31) Chrysene	21.92	228	1916m	0.022	ng	

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

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