

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE120517\
 Data File : BE094936.D
 Acq On : 6 Dec 2017 3:58
 Operator : SJ/JU
 Sample : I6698-14
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 FT-DUP03-20171130

Manual Integrations
 APPROVED

Sohil
 12/6/2017 5:10:41 PM

Quant Time: Dec 06 07:04:39 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	8651	0.40	ng	0.00
7) Naphthalene-d8	11.33	136	18527	0.40	ng	0.00
13) Acenaphthene-d10	15.06	164	10808	0.40	ng	-0.02
19) Phenanthrene-d10	17.77	188	26499	0.40	ng	0.00
27) Chrysene-d12	21.87	240	28195	0.40	ng	-0.01
34) Perylene-d12	24.55	264	23331	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.95	112	2577	0.19	ng	0.00
5) Phenol-d6	7.60	99	2340	0.12	ng	0.01
8) Nitrobenzene-d5	9.67	82	5744	0.52	ng	0.00
11) 2-Methylnaphthalene-d10	12.87	152	11919	0.40	ng	0.00
14) 2,4,6-Tribromophenol	16.53	330	1315	0.55	ng	0.00
15) 2-Fluorobiphenyl	13.71	172	18644	0.40	ng	0.00
25) Fluoranthene-d10	19.75	212	142694	0.39	ng	0.00
29) Terphenyl-d14	20.31	244	28676	0.55	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	3.67	88	992	0.110	ng	Qvalue # 53
30) Benzo(a)anthracene	21.86	228	3194	0.037	ng	# 66
31) Chrysene	21.92	228	2974m	0.031	ng	
32) Bis(2-ethylhexyl)phthalate	21.75	149	23377	0.163	ng	100
33) Indeno(1,2,3-cd)pyrene	27.43	276	3038	0.039	ng	95
35) Benzo(b)fluoranthene	23.72	252	3259	0.038	ng	# 63
36) Benzo(k)fluoranthene	23.77	252	2735	0.031	ng	# 59
37) Benzo(a)pyrene	24.43	252	2757	0.037	ng	# 71
38) Dibenzo(a,h)anthracene	27.44	278	2527	0.037	ng	# 68
39) Benzo(a,h,i)perylene	28.30	276	2810	0.041	ng	# 61

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

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