

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE040121\
 Data File : BE101178.D
 Acq On : 1 Apr 2021 15:29
 Operator : CG/JU
 Sample : M1843-17
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 DUP08-20210325

Quant Time: Apr 01 16:13:01 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-SIM-BE032421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 01 12:49:55 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.842	152	4445	0.40	ng	0.00
7) Naphthalene-d8	10.627	136	18904	0.40	ng	# 0.00
13) Acenaphthene-d10	14.457	164	12048	0.40	ng	0.00
19) Phenanthrene-d10	17.211	188	30531	0.40	ng	# 0.00
29) Chrysene-d12	21.366	240	35756	0.40	ng	# 0.00
36) Perylene-d12	23.848	264	32408	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.435	112	2038	0.19	ng	0.00
5) Phenol-d6	7.001	99	1830	0.11	ng	0.00
8) Nitrobenzene-d5	8.977	82	5584	0.49	ng	0.00
11) 2-Methylnaphthalene-d10	12.209	152	12137	0.39	ng	0.00
14) 2,4,6-Tribromophenol	15.956	330	1315	0.61	ng	0.00
15) 2-Fluorobiphenyl	13.088	172	17187	0.38	ng	0.00
27) Fluoranthene-d10	19.227	212	171582	0.44	ng	0.00
31) Terphenyl-d14	19.824	244	39655	0.49	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.373	88	31297	4.34	ng	95
6) bis(2-Chloroethyl)ether	7.266	93	940	0.07	ng	95
25) Phenanthrene	17.257	178	1964	0.02	ng	# 96
34) Bis(2-ethylhexyl)phtha...	21.293	149	9815	0.16	ng	100
40) Dibenzo(a,h)anthracene	26.491	278	521	0.11	ng	# 35

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

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