

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100421\
 Data File : BN016746.D
 Acq On : 05 Oct 2021 14:28
 Operator : CG/JU
 Sample : M3829-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE105D2-20210917

Quant Time: Oct 05 16:00:19 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 04 17:52:33 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.633	152	25033	0.400	ng	0.00
7) Naphthalene-d8	10.407	136	115196	0.400	ng	# 0.00
13) Acenaphthene-d10	14.269	164	61808	0.400	ng	0.00
19) Phenanthrene-d10	17.030	188	120423	0.400	ng	0.00
29) Chrysene-d12	21.238	240	118052	0.400	ng	0.00
35) Perylene-d12	23.491	264	120086	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.264	112	6903	0.108	ng	0.00
5) Phenol-d6	6.822	99	4245	0.060	ng	0.00
8) Nitrobenzene-d5	8.785	82	20870	0.258	ng	0.00
11) 2-Methylnaphthalene-d10	12.003	152	50953	0.297	ng	0.00
14) 2,4,6-Tribromophenol	15.779	330	8845	0.365	ng	0.00
15) 2-Fluorobiphenyl	12.899	172	71127	0.288	ng	0.00
27) Fluoranthene-d10	19.068	212	133558	0.405	ng	0.00
31) Terphenyl-d14	19.679	244	114025	0.432	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.253	88	45881	3.927	ng	# 68
24) Pentachlorophenol	16.689	266	126	0.342	ng	91
34) Bis(2-ethylhexyl)phtha...	21.174	149	34636	0.156	ng	99

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

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