

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN113023\  
 Data File : BN028948.D  
 Acq On : 01 Dec 2023 15:29  
 Operator : MA/JU  
 Sample : 05350-10  
 Misc :  
 ALS Vial : 39 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 RE115D2-20231107

Quant Time: Dec 01 16:20:42 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN113023.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Dec 01 00:28:34 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.797	152	4421	0.400	ng	-0.01	
7) Naphthalene-d8	10.584	136	11643	0.400	ng	-0.01	
13) Acenaphthene-d10	14.428	164	4844	0.400	ng	0.00	
19) Phenanthrene-d10	17.171	188	7593	0.400	ng	-0.01	
29) Chrysene-d12	21.374	240	5005	0.400	ng	# 0.00	
35) Perylene-d12	23.708	264	4920	0.400	ng	# 0.00	
System Monitoring Compounds							
4) 2-Fluorophenol	5.420	112	2165	0.162	ng	-0.03	
5) Phenol-d6	7.017	99	1428	0.096	ng	-0.03	
8) Nitrobenzene-d5	8.971	82	3386	0.293	ng	-0.01	
11) 2-Methylnaphthalene-d10	12.174	152	5352	0.286	ng	-0.01	
14) 2,4,6-Tribromophenol	15.918	330	626	0.191	ng	-0.01	
15) 2-Fluorobiphenyl	13.060	172	9435	0.451	ng	-0.01	
27) Fluoranthene-d10	19.209	212	7321	0.367	ng	0.00	
31) Terphenyl-d14	19.817	244	6018	0.414	ng	0.00	
Target Compounds							
2) 1,4-Dioxane	3.340	88	29300	6.767	ng		95
6) bis(2-Chloroethyl)ether	7.255	93	546	0.044	ng		91
16) Acenaphthylene	14.140	152	6254	0.244	ng		96
17) Acenaphthene	14.482	154	756	0.044	ng		94
18) Fluorene	15.497	166	852	0.040	ng	#	11
30) Pyrene	19.604	202	767	0.114	ng	#	80
32) Benzo(a)anthracene	21.356	228	552	0.026	ng	#	1
33) Chrysene	21.409	228	736	0.035	ng	#	1

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

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