

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG111222\  
 Data File : BG055600.D  
 Acq On : 12 Nov 2022 22:15  
 Operator : CG/JU  
 Sample : N5575-01 2X  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 OK-1-111022

Manual Integrations  
 APPROVED

Reviewed By :Christian Giraldo 11/14/2022  
 Supervised By :Jagrut Upadhyay 11/14/2022

Quant Time: Nov 14 03:32:47 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG110922.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Nov 10 03:39:46 2022  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
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Internal Standards							
1) 1,4-Dichlorobenzene-d4	8.270	152	90808	20.000	ng	0.00	
21) Naphthalene-d8	11.102	136	370849	20.000	ng	0.00	
39) Acenaphthene-d10	14.892	164	240416	20.000	ng	0.00	
64) Phenanthrene-d10	17.630	188	495267	20.000	ng	0.00	
76) Chrysene-d12	21.936	240	460430	20.000	ng	0.00	
86) Perylene-d12	25.373	264	389022	20.000	ng	0.02	
System Monitoring Compounds							
5) 2-Fluorophenol	5.796	112	121660	25.519	ng	0.00	
7) Phenol-d6	7.406	99	159891	24.279	ng	0.00	
23) Nitrobenzene-d5	9.439	82	101136	18.488	ng	0.00	
42) 2,4,6-Tribromophenol	16.372	330	70332	29.983	ng	0.00	
45) 2-Fluorobiphenyl	13.517	172	250442	15.713	ng	0.00	
79) Terphenyl-d14	20.215	244	360201	15.713	ng	0.00	
Target Compounds							
71) Phenanthrene	17.677	178	113499	4.264	ng		Qvalue 98
75) Fluoranthene	19.668	202	173254	5.458	ng		99
78) Pyrene	20.033	202	144740	4.645	ng		96
81) Benzo(a)anthracene	21.913	228	67726	2.147	ng	#	92
83) Chrysene	21.983	228	75934m	2.532	ng		
88) Benzo(b)fluoranthene	24.275	252	74516	3.023	ng	#	68
90) Benzo(a)pyrene	25.203	252	55032	2.592	ng	#	73
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

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