

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN040523\  
 Data File : BN024760.D  
 Acq On : 05 Apr 2023 00:08  
 Operator : CG/JU  
 Sample : PB151661BL  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

**Instrument :**  
 BNA\_N  
**ClientSampleId :**  
 PB151661BL

Quant Time: Apr 05 02:30:46 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN040523.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Apr 05 02:15:45 2023  
 Response via : Initial Calibration

**Manual Integrations**  
**APPROVED**  
 Reviewed By :Christian Giraldo 04/06/2023  
 Supervised By :Jagrut Upadhyay 04/06/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.960	152	9679	0.400	ng	0.00
7) Naphthalene-d8	10.770	136	25459	0.400	ng	0.00
13) Acenaphthene-d10	14.600	164	14292	0.400	ng	0.00
19) Phenanthrene-d10	17.338	188	31826	0.400	ng	0.00
29) Chrysene-d12	21.529	240	28415	0.400	ng	0.00
35) Perylene-d12	23.968	264	26396	0.400	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.505	112	8755	0.392	ng	0.00
5) Phenol-d6	7.115	99	10399	0.369	ng	0.00
8) Nitrobenzene-d5	9.126	82	7713	0.376	ng	0.00
11) 2-Methylnaphthalene-d10	12.357	152	13951	0.414	ng	0.00
14) 2,4,6-Tribromophenol	16.122	330	1364m	0.212	ng	0.04
15) 2-Fluorobiphenyl	13.221	172	20880	0.382	ng	0.00
27) Fluoranthene-d10	19.370	212	29570	0.410	ng	0.00
31) Terphenyl-d14	19.964	244	26258	0.408	ng	0.00

Target Compounds Qvalue

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

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