

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110221\
 Data File : BM032855.D
 Acq On : 02 Nov 2021 17:19
 Operator : CG/JU
 Sample : PB140166BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB140166BL

Quant Time: Nov 02 17:54:58 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM110221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 02 16:07:18 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.489	152	118511	20.000	ng	0.00
21) Naphthalene-d8	10.272	136	509229	20.000	ng	# 0.00
39) Acenaphthene-d10	14.154	164	361406	20.000	ng	0.00
64) Phenanthrene-d10	16.889	188	794171	20.000	ng	0.00
76) Chrysene-d12	21.071	240	735079	20.000	ng	# 0.00
86) Perylene-d12	23.188	264	692995	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.084	112	1172897	172.542	ng	0.00
7) Phenol-d6	6.695	99	1531778	151.314	ng	0.00
23) Nitrobenzene-d5	8.648	82	955235	102.670	ng	0.00
42) 2,4,6-Tribromophenol	15.648	330	583719	144.217	ng	0.00
45) 2-Fluorobiphenyl	12.766	172	2255050	96.776	ng	0.00
79) Terphenyl-d14	19.524	244	3268455	97.076	ng	0.00

Target Compounds	Qvalue

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

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