

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080221\
 Data File : BP006465.D
 Acq On : 03 Aug 2021 00:37
 Operator : CG/JU
 Sample : PB138104BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB138104BL

Quant Time: Aug 03 05:39:55 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 03 05:37:51 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.757	152	245429	20.000	ng	0.00
21) Naphthalene-d8	10.540	136	1027073	20.000	ng	# 0.00
39) Acenaphthene-d10	14.387	164	560388	20.000	ng	0.00
64) Phenanthrene-d10	17.139	188	1119142	20.000	ng	# 0.00
76) Chrysene-d12	21.239	240	1067453	20.000	ng	# 0.00
86) Perylene-d12	23.568	264	1121061	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.358	112	1793396	113.954	ng	0.00
7) Phenol-d6	6.940	99	2259508	101.693	ng	0.00
23) Nitrobenzene-d5	8.916	82	1494462	68.601	ng	0.00
42) 2,4,6-Tribromophenol	15.881	330	580092	101.236	ng	0.00
45) 2-Fluorobiphenyl	13.016	172	2548643	68.170	ng	0.00
79) Terphenyl-d14	19.716	244	3701260	69.818	ng	0.00

Target Compounds	Qvalue

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

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