

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122721\
 Data File : BP008561.D
 Acq On : 27 Dec 2021 18:11
 Operator : CG/JU
 Sample : PB141663BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK663

Quant Time: Dec 28 01:30:24 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 23 14:48:11 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	541168	20.000	ng/u1	0.00
20) Naphthalene-d8	10.704	136	2221463	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.534	164	1527035	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.280	188	3189929	20.000	ng/u1	0.00
79) Chrysene-d12	21.363	240	2531977	20.000	ng/u1	0.00
88) Perylene-d12	23.792	264	2089046	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	76371	6.361	ng/uL	0.00
4) Pyridine-d5	3.758	84	986462	29.360	ng/u1	0.00
7) Phenol-d5	7.057	99	1301044	30.136	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	741040	30.221	ng/u1	0.00
11) 2-Chlorophenol-d4	7.434	132	1072200	31.220	ng/u1	0.00
15) 4-Methylphenol-d8	8.604	113	1047949	30.259	ng/u1	0.00
21) Nitrobenzene-d5	9.057	128	513185	32.647	ng/u1	0.00
24) 2-Nitrophenol-d4	9.781	143	487752	33.841	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.316	165	1084904	30.664	ng/u1	0.00
31) 4-Chloroaniline-d4	10.834	131	1563534	32.015	ng/u1	0.00
46) Dimethylphthalate-d6	13.939	166	3756752	33.297	ng/u1	0.00
49) Acenaphthylene-d8	14.222	160	4558361	32.311	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	497833	27.836	ng/u1	0.00
60) Fluorene-d10	15.522	176	3241542	33.091	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.633	200	338043	24.074	ng/u1	0.00
73) Anthracene-d10	17.380	188	5091724	33.684	ng/u1	0.00
81) Pyrene-d10	19.610	212	5561096	36.654	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.633	264	3619314	33.662	ng/u1	0.00

Target Compounds Qvalue

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

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