

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP012007.D
 Acq On : 07 Oct 2022 13:47
 Operator : CG/JU
 Sample : PB148168BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK168

Quant Time: Oct 07 23:08:04 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 07 03:22:44 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	202584	20.000	ng/u1	0.00
20) Naphthalene-d8	10.681	136	839645	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.522	164	491260	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.281	188	1020602	20.000	ng/u1	0.00
79) Chrysene-d12	21.369	240	987922	20.000	ng/u1	0.00
88) Perylene-d12	23.804	264	1009266	20.000	ng/u1	0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	25769	5.610	ng/uL	0.00
4) Pyridine-d5	3.723	84	322529	23.971	ng/u1	0.00
7) Phenol-d5	7.040	99	487861	27.895	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	277709	28.647	ng/u1	0.00
11) 2-Chlorophenol-d4	7.405	132	407803	29.538	ng/u1	0.00
15) 4-Methylphenol-d8	8.587	113	382444	26.453	ng/u1	0.00
21) Nitrobenzene-d5	9.046	128	199840	31.247	ng/u1	0.00
24) 2-Nitrophenol-d4	9.769	143	214910	31.408	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.305	165	355735	28.108	ng/u1	0.00
31) 4-Chloroaniline-d4	10.828	131	509900	26.427	ng/u1	0.00
46) Dimethylphthalate-d6	13.934	166	980610	27.914	ng/u1	0.00
49) Acenaphthylene-d8	14.216	160	1251444	31.248	ng/u1	0.00
54) 4-Nitrophenol-d4	14.734	143	163030	22.410	ng/u1	0.01
60) Fluorene-d10	15.516	176	956831	31.363	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	103347	16.730	ng/u1	0.00
73) Anthracene-d10	17.381	188	1529766	32.905	ng/u1	0.00
81) Pyrene-d10	19.616	212	1726806	34.514	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.645	264	1717517	33.995	ng/u1	0.01

Target Compounds	Qvalue

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

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