

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012117.D
 Acq On : 13 Oct 2022 13:00
 Operator : CG/JU
 Sample : PB148253BL
 Misc :
 ALS Vial : 88 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 13 23:33:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 12 23:03:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	270967	20.000	ng/ul	0.00
20) Naphthalene-d8	10.657	136	1092200	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.498	164	595986	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.257	188	1224680	20.000	ng/ul	0.00
79) Chrysene-d12	21.345	240	1304054	20.000	ng/ul	0.00
88) Perylene-d12	23.769	264	1364899	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	35071	5.103	ng/uL	0.00
4) Pyridine-d5	3.699	84	449008	24.566	ng/ul	0.00
7) Phenol-d5	7.016	99	600745	27.566	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	339559	27.450	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	499800	28.893	ng/ul	0.00
15) 4-Methylphenol-d8	8.563	113	456564	27.276	ng/ul	0.00
21) Nitrobenzene-d5	9.016	128	237654	30.809	ng/ul	0.00
24) 2-Nitrophenol-d4	9.740	143	243477	31.129	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	418184	27.588	ng/ul	0.00
31) 4-Chloroaniline-d4	10.799	131	632627	28.321	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	1131533	29.979	ng/ul	0.00
49) Acenaphthylene-d8	14.193	160	1377521	29.513	ng/ul	0.00
54) 4-Nitrophenol-d4	14.710	143	195905	26.437	ng/ul	0.00
60) Fluorene-d10	15.492	176	1032516	30.235	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.616	200	131583	20.348	ng/ul	0.00
73) Anthracene-d10	17.357	188	1629838	30.647	ng/ul	0.00
81) Pyrene-d10	19.592	212	1965455	29.420	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.610	264	2049041	30.882	ng/ul	0.00

Target Compounds Qvalue

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

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