

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011904.D
 Acq On : 04 Oct 2022 10:26
 Operator : CG/JU
 Sample : PB148002BL
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK002

Quant Time: Oct 05 04:29:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 01 00:53:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	274781	20.000	ng/u1	0.00
20) Naphthalene-d8	10.693	136	1053524	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.528	164	600869	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.292	188	1246856	20.000	ng/u1	0.00
79) Chrysene-d12	21.375	240	1347355	20.000	ng/u1	-0.01
88) Perylene-d12	23.810	264	1399735	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	44117	7.081	ng/uL	0.00
4) Pyridine-d5	3.723	84	531546	29.126	ng/u1	-0.01
7) Phenol-d5	7.046	99	739532	31.175	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.217	67	412989	31.409	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.417	132	624685	33.359	ng/u1	0.00
15) 4-Methylphenol-d8	8.593	113	574340	29.289	ng/u1	-0.01
21) Nitrobenzene-d5	9.052	128	285852	35.622	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.775	143	300765	35.032	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.316	165	535456	33.719	ng/u1	0.00
31) 4-Chloroaniline-d4	10.834	131	779840	32.212	ng/u1	-0.01
46) Dimethylphthalate-d6	13.946	166	1618343	37.664	ng/u1	0.00
49) Acenaphthylene-d8	14.222	160	1885158	38.485	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.734	143	295119	33.167	ng/u1	0.00
60) Fluorene-d10	15.528	176	1425653	38.205	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	227733	30.176	ng/u1	0.00
73) Anthracene-d10	17.387	188	2296182	40.428	ng/u1	-0.01
81) Pyrene-d10	19.622	212	2829032	41.460	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.651	264	2982466	42.565	ng/u1	-0.01

Target Compounds Qvalue

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

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