

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011989.D
 Acq On : 07 Oct 2022 02:37
 Operator : CG/JU
 Sample : PB148077BL
 Misc :
 ALS Vial : 78 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK077

Quant Time: Oct 07 03:26:45 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	258772	20.000	ng/u1	0.00
20) Naphthalene-d8	10.681	136	1098767	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.522	164	686663	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.275	188	1456795	20.000	ng/u1	0.00
79) Chrysene-d12	21.363	240	1354153	20.000	ng/u1	0.00
88) Perylene-d12	23.792	264	1367041	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	36019	6.139	ng/uL	0.00
4) Pyridine-d5	3.716	84	438290	25.502	ng/u1	0.00
7) Phenol-d5	7.040	99	635177	28.433	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	364419	29.429	ng/u1	0.00
11) 2-Chlorophenol-d4	7.404	132	522043	29.602	ng/u1	0.00
15) 4-Methylphenol-d8	8.587	113	499492	27.047	ng/u1	0.00
21) Nitrobenzene-d5	9.040	128	253356	30.273	ng/u1	0.00
24) 2-Nitrophenol-d4	9.763	143	265914	29.697	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.304	165	473715	28.602	ng/u1	0.00
31) 4-Chloroaniline-d4	10.822	131	691837	27.400	ng/u1	0.00
46) Dimethylphthalate-d6	13.934	166	1477167	30.083	ng/u1	0.00
49) Acenaphthylene-d8	14.216	160	1702936	30.421	ng/u1	0.00
54) 4-Nitrophenol-d4	14.722	143	241623	23.762	ng/u1	0.00
60) Fluorene-d10	15.516	176	1328511	31.154	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.633	200	184307	20.902	ng/u1	0.00
73) Anthracene-d10	17.375	188	2155579	32.483	ng/u1	0.00
81) Pyrene-d10	19.616	212	2447121	35.683	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.633	264	2316583	33.852	ng/u1	0.00

Target Compounds Qvalue

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

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