

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011914.D
 Acq On : 04 Oct 2022 16:08
 Operator : CG/JU
 Sample : N4873-16
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXZ61

Quant Time: Oct 05 05:24:20 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	252224	20.000	ng/u1	0.00
20) Naphthalene-d8	10.693	136	1041367	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.534	164	635831	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.286	188	1380983	20.000	ng/u1	-0.01
79) Chrysene-d12	21.374	240	1477525	20.000	ng/u1	-0.01
88) Perylene-d12	23.804	264	1484731	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	28492	4.982	ng/uL	0.00
4) Pyridine-d5	3.728	84	100207	5.982	ng/u1	0.00
7) Phenol-d5	7.052	99	123446	5.669	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	352360	29.194	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.416	132	391640	22.784	ng/u1	0.00
15) 4-Methylphenol-d8	8.593	113	234195	13.011	ng/u1	-0.01
21) Nitrobenzene-d5	9.052	128	262371	33.078	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.775	143	249588	29.411	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.310	165	428359	27.290	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.834	131	420037	17.553	ng/u1	-0.01
46) Dimethylphthalate-d6	13.940	166	1474393	32.427	ng/u1	-0.01
49) Acenaphthylene-d8	14.222	160	1689128	32.587	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.734	143	57328	6.089	ng/u1	0.00
60) Fluorene-d10	15.528	176	1337839	33.881	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	217129	25.977	ng/u1	0.00
73) Anthracene-d10	17.386	188	2197550	34.934	ng/u1	-0.01
81) Pyrene-d10	19.622	212	2769901	37.017	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.651	264	2723172	36.639	ng/u1	-0.01

Target Compounds Qvalue

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

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