

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013200.D
 Acq On : 21 Dec 2022 10:40
 Operator : CG/JU
 Sample : SP5089
 Misc : SFAM SURROGATE
 ALS Vial : 82 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SP5089

Quant Time: Dec 21 21:51:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 19 21:43:11 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.951	152	498400	20.000	ng/u1	0.00
20) Naphthalene-d8	10.763	136	1953533	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.592	164	1095443	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.351	188	2133219	20.000	ng/u1	0.00
79) Chrysene-d12	21.433	240	2178974	20.000	ng/u1	0.00
88) Perylene-d12	23.915	264	2560406	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	197070	16.906	ng/uL	0.00
4) Pyridine-d5	3.757	84	2905313	87.740	ng/u1	0.00
7) Phenol-d5	7.110	99	3383367	83.343	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	2060972	84.159	ng/u1	0.00
11) 2-Chlorophenol-d4	7.481	132	2742939	85.831	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	2605589	78.333	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	1360695	94.801	ng/u1	0.00
24) 2-Nitrophenol-d4	9.845	143	1553264	96.125	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	2633267	83.899	ng/u1	0.00
31) 4-Chloroaniline-d4	10.904	131	3503422	79.068	ng/u1	0.00
46) Dimethylphthalate-d6	14.004	166	6654597	79.043	ng/u1	0.00
49) Acenaphthylene-d8	14.286	160	8048473	87.001	ng/u1	0.00
54) 4-Nitrophenol-d4	14.792	143	1097121	72.446	ng/u1	0.00
60) Fluorene-d10	15.586	176	5715325	80.243	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.710	200	1002260	73.010	ng/u1	0.00
73) Anthracene-d10	17.451	188	8293328	86.174	ng/u1	0.00
81) Pyrene-d10	19.680	212	9823509	78.542	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.762	264	11397880	89.310	ng/u1	0.00

Target Compounds Qvalue

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

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