

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122722\
 Data File : BP013343.D
 Acq On : 27 Dec 2022 17:43
 Operator : CG/JU
 Sample : PB149814BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK814

Quant Time: Dec 28 00:20:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 28 00:16:35 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	569378	20.000	ng/u1	0.00
20) Naphthalene-d8	10.746	136	2098428	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.575	164	1283291	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.334	188	2999474	20.000	ng/u1	0.00
79) Chrysene-d12	21.422	240	1988457	20.000	ng/u1	0.00
88) Perylene-d12	23.886	264	1932801	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.323	96	410	0.031	ng/uL	0.00
4) Pyridine-d5	3.746	84	61	0.002	ng/u1	0.00
7) Phenol-d5	7.093	99	7	0.000	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	418	0.015	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.464	132	37	0.001	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	22	0.001	ng/u1	-0.01
21) Nitrobenzene-d5	9.093	128	65	0.004	ng/u1	0.00
24) 2-Nitrophenol-d4	9.828	143	8	0.000	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	105	0.003	ng/u1	0.01
31) 4-Chloroaniline-d4	10.887	131	161	0.003	ng/u1	0.00
46) Dimethylphthalate-d6	13.987	166	12	0.000	ng/u1	0.00
49) Acenaphthylene-d8	14.257	160	49	0.000	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.787	143	30	0.002	ng/u1	0.00
60) Fluorene-d10	15.575	176	1336	0.016	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	8	0.000	ng/u1	0.00
73) Anthracene-d10	17.334	188	2999474	22.166	ng/u1	-0.10
81) Pyrene-d10	19.669	212	6722	0.059	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.886	264	1726138	17.917	ng/u1	0.15

Target Compounds Qvalue

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

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