

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082823\
 Data File : BP016840.D
 Acq On : 29 Aug 2023 19:38
 Operator : MA/JU
 Sample : 04134-09
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCHK2

Quant Time: Aug 29 23:22:27 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 29 01:12:28 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	351557	20.000	ng/u1	0.00
20) Naphthalene-d8	10.858	136	1568669	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.681	164	995267	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.445	188	2138088	20.000	ng/u1	-0.01
79) Chrysene-d12	21.545	240	1510310	20.000	ng/u1	0.00
88) Perylene-d12	24.110	264	1367106	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	43721	5.140	ng/uL	0.00
4) Pyridine-d5	3.817	84	252799	9.725	ng/u1	0.00
7) Phenol-d5	7.169	99	251889	8.058	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.364	67	713021	35.787	ng/u1	0.00
11) 2-Chlorophenol-d4	7.546	132	657516	27.975	ng/u1	0.00
15) 4-Methylphenol-d8	8.728	113	461668	18.091	ng/u1	0.00
21) Nitrobenzene-d5	9.222	128	426832	35.987	ng/u1	0.00
24) 2-Nitrophenol-d4	9.940	143	412239	32.974	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.463	165	762774	32.727	ng/u1	0.00
31) 4-Chloroaniline-d4	11.010	131	968460	27.095	ng/u1	0.00
46) Dimethylphthalate-d6	14.093	166	2835152	37.773	ng/u1	0.00
49) Acenaphthylene-d8	14.381	160	3226504	38.026	ng/u1	0.00
54) 4-Nitrophenol-d4	14.887	143	73426	4.733	ng/u1	0.00
60) Fluorene-d10	15.675	176	2530803	40.040	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200	326635	25.653	ng/u1	0.00
73) Anthracene-d10	17.551	188	4454099	46.998	ng/u1	0.00
81) Pyrene-d10	19.781	212	4831801	58.250	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.945	264	3252054	47.489	ng/u1	0.00

Target Compounds Qvalue

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

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